

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020624\
 Data File : BN029531.D
 Acq On : 06 Feb 2024 11:07
 Operator : MA/JU
 Sample : SSTD0.251
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.2224

Quant Time: Feb 07 02:11:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 02:09:40 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	4620	0.400	ng/ul	0.00
4) Naphthalene-d8	10.689	136	12365	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.530	164	5395	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.284	188	9589	0.400	ng/ul	0.00
17) Chrysene-d12	21.481	240	6381	0.400	ng/ul	0.00
23) Perylene-d12	23.858	264	6477	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	1189	0.233	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.283	152	3115	0.167	ng/ul	0.00
18) Fluoranthene-d10	19.318	212	4429	0.216	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.374	88	1302	0.231	ng/ul#	90
5) Naphthalene	10.738	128	7093	0.205	ng/ul	99
7) 2-Methylnaphthalene	12.355	142	3948	0.173	ng/ul	99
8) 1-Methylnaphthalene	12.575	142	4061	0.175	ng/ul	99
10) Acenaphthylene	14.252	152	5370	0.199	ng/ul	98
11) Acenaphthene	14.594	153	4101	0.207	ng/ul	98
12) Fluorene	15.584	166	4169	0.181	ng/ul	99
14) Pentachlorophenol	16.930	266	408	0.177	ng/ul#	49
15) Phenanthrene	17.327	178	6130	0.208	ng/ul	98
16) Anthracene	17.415	178	5000	0.185	ng/ul	98
19) Fluoranthene	19.346	202	6066	0.224	ng/ul	97
20) Pyrene	19.708	202	6361	0.226	ng/ul	97
21) Benzo(a)anthracene	21.467	228	4601	0.182	ng/ul	99
22) Chrysene	21.519	228	5201	0.211	ng/ul	99
24) Benzo(b)fluoranthene	23.136	252	4847	0.199	ng/ul	89
25) Benzo(k)fluoranthene	23.185	252	5585	0.219	ng/ul#	87
26) Benzo(a)pyrene	23.753	252	4760	0.207	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	26.322	276	5738	0.225	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.352	278	4464	0.219	ng/ul#	86
29) Benzo(g,h,i)perylene	27.067	276	5361	0.257	ng/ul	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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